

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072622\
 Data File : BP011103.D
 Acq On : 26 Jul 2022 11:02
 Operator : CG/JU
 Sample : PB146404BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK404

Quant Time: Jul 26 14:40:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 26 14:31:36 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	430589	20.000	ng/u1	0.00
20) Naphthalene-d8	10.457	136	1978658	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.328	164	1206235	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.092	188	2245668	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.210	240	1432851	20.000	ng/u1	# 0.00
88) Perylene-d12	23.557	264	1314534	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	77271	6.470	ng/uL	0.00
4) Pyridine-d5	3.569	84	1062510	34.417	ng/u1	0.00
7) Phenol-d5	6.840	99	1255519	35.521	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	868206	38.199	ng/u1	0.00
11) 2-Chlorophenol-d4	7.199	132	1045645	36.612	ng/u1	0.00
15) 4-Methylphenol-d8	8.381	113	1066358	37.685	ng/u1	0.00
21) Nitrobenzene-d5	8.828	128	545262	37.814	ng/u1	0.00
24) 2-Nitrophenol-d4	9.546	143	569957	39.946	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	916356	35.118	ng/u1	0.00
31) 4-Chloroaniline-d4	10.604	131	1536698	35.723	ng/u1	0.00
46) Dimethylphthalate-d6	13.751	166	3379946	40.998	ng/u1	0.00
49) Acenaphthylene-d8	14.016	160	3856743	39.649	ng/u1	0.00
54) 4-Nitrophenol-d4	14.545	143	471190	28.240	ng/u1	0.00
60) Fluorene-d10	15.328	176	2645212	37.748	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	394549	35.506	ng/u1	0.00
73) Anthracene-d10	17.192	188	3769620	39.485	ng/u1	0.00
81) Pyrene-d10	19.451	212	3643438	48.416	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.404	264	2568705	40.685	ng/u1	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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